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* * * * * * * * * *   Welcome to STN International   * * * * * * * * * *

 NEWS  1          Web Page for STN Seminar Schedule - N. America
 NEWS  2  DEC 01  ChemPort single article sales feature unavailable
 NEWS  3  APR 03  CAS coverage of exemplified prophetic substances
                  enhanced
 NEWS  4  APR 07  STN is raising the limits on saved answers
 NEWS  5  APR 24  CA/CAplus now has more comprehensive patent assignee
                  information
 NEWS  6  APR 26  USPATFULL and USPAT2 enhanced with patent
                  assignment/reassignment information
 NEWS  7  APR 28  CAS patent authority coverage expanded
 NEWS  8  APR 28  ENCOMPLIT/ENCOMPLIT2 search fields enhanced
 NEWS  9  APR 28  Limits doubled for structure searching in CAS
                  REGISTRY
 NEWS 10  MAY 08  STN Express, Version 8.4, now available
 NEWS 11  MAY 11  STN on the Web enhanced
 NEWS 12  MAY 11  BEILSTEIN substance information now available on
                  STN Easy
 NEWS 13  MAY 14  DGENE, PCTGEN and USGENE enhanced with increased
                  limits for exact sequence match searches and
                  introduction of free HIT display format
 NEWS 14  MAY 15  INPADOCDB and INPAFAMDB enhanced with Chinese legal
                  status data
 NEWS 15  MAY 28  CAS databases on STN enhanced with NANO super role in
                  records back to 1992
 NEWS 16  JUN 01  CAS REGISTRY Source of Registration (SR) searching
                  enhanced on STN
 NEWS 17  JUN 26  NUTRACEUT and PHARMAML no longer updated

 NEWS 18  JUN 29  IMSCOPROFILE now reloaded monthly

 NEWS 19  JUN 29  EPFULL adds SLART to AB, MCLM, and TI fields

 NEWS EXPRESS MAY 26 09 CURRENT WINDOWS VERSION IS V8.4,
              AND CURRENT DISCOVER FILE IS DATED 06 APRIL 2009.

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FILE 'HOME' ENTERED AT 11:33:37 ON 01 JUL 2009

=> file reg
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                       ENTRY    SESSION
FULL ESTIMATED COST                                     0.22       0.22

FILE 'REGISTRY' ENTERED AT 11:33:57 ON 01 JUL 2009
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STRUCTURE FILE UPDATES: 30 JUN 2009 HIGHEST RN 1160555-05-4
DICTIONARY FILE UPDATES: 30 JUN 2009 HIGHEST RN 1160555-05-4

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=>
Uploading C:\Program Files\Stnexp\Queries\11573868-Ib.str

L1   STRUCTURE UPLOADED

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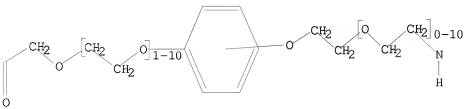
L1 HAS NO ANSWERS
L1   STR



Structure attributes must be viewed using STN Express query preparation.

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SAMPLE SCREEN SEARCH COMPLETED -   5144 TO ITERATE

 38.9% PROCESSED   2000 ITERATIONS   0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS: 98579 TO   107181
PROJECTED ANSWERS: 0 TO   0

L2   0 SEA SSS SAM L1

=> s l1 full
FULL SEARCH INITIATED 11:34:25 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -   101816 TO ITERATE

100.0% PROCESSED   101816 ITERATIONS   6 ANSWERS
SEARCH TIME: 00.00.01

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=> d l3 scan

L3    6 ANSWERS   REGISTRY  COPYRIGHT 2009 ACS on STN
IN    5,8,11,14-Tetraoxa-2-azahexadecanoic acid,
      16-[3-[(16-carboxy-3,6,9,12,15-pentaoxahexadec-1-yl)oxy]-5-[[18-(9H-
      fluoren-9-yl)-16-oxo-3,6,9,12,17-pentaoxa-15-azaoctadec-1-yl]oxy]phenoxy]-
      , 1-(9H-fluoren-9-ylmethyl) ester
MF    C68 H90 N2 O22
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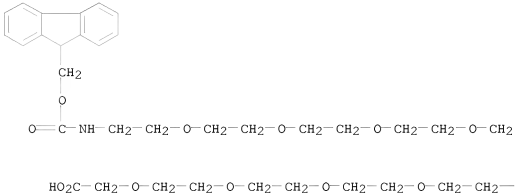
PAGE 1-A



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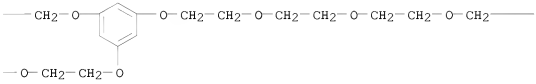

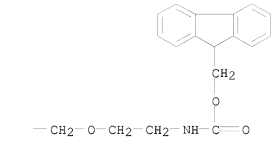

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):file caplus
'FILE CAPLUS' IS NOT VALID HERE

To display more answers, enter the number of answers you would like to see. To end the display, enter "NONE", "N", "0", or "END".
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> file caplus

COST IN U.S. DOLLARS	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	186.36	186.58

FILE 'CAPLUS' ENTERED AT 11:34:52 ON 01 JUL 2009
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FILE COVERS 1907 - 1 Jul 2009 VOL 151 ISS 1
FILE LAST UPDATED: 30 Jun 2009 (20090630/ED)
REVISED CLASS FIELDS (/NCL) LAST RELOADED: Apr 2009
USPTO MANUAL OF CLASSIFICATIONS THESAURUS ISSUE DATE: Apr 2009

CAplus now includes complete International Patent Classification (IPC) reclassification data for the second quarter of 2009.

CAS Information Use Policies apply and are available at:

http://www.cas.org/legal/infopolicy.html

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3
L4 3 L3

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=> d l4 ibib abs hitstr 1-
YOU HAVE REQUESTED DATA FROM 3 ANSWERS - CONTINUE? Y/(N):y

L4   ANSWER 1 OF 3  CAPLUS  COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER:         2004:406941  CAPLUS
DOCUMENT NUMBER:          141:273853
TITLE:                    Design and synthesis of novel hydrophilic spacers for
                          the reduction of nonspecific binding proteins on
                          affinity resins
AUTHOR(S):                Shiyama, Takaaki; Furuya, Minoru; Yamazaki, Akira;
                          Terada, Tomohiro; Tanaka, Akito
CORPORATE SOURCE:         Chemistry Department, Reverse Proteomics Research
                          Institute Co., Ltd, Chiba, 292-0818, Japan
SOURCE:                   Bioorganic & Medicinal Chemistry (2004), 12(11),
                          2831-2841
                          CODEN: BMECEP; ISSN: 0968-0896
PUBLISHER:                Elsevier Ltd.
DOCUMENT TYPE:            Journal
LANGUAGE:                 English
OTHER SOURCE(S):          CASREACT 141:273853
AB   Tubulin and actin often bind nonspecifically to affinity chromatog.
     resins, complicating research toward identifying the cellular targets.
     Reduction of nonspecific binding proteins is important for success in finding
     such targets.  We herein disclose the design, synthesis, and effectiveness
     in reduction of nonspecific binding proteins, of novel hydrophilic spacers
     (2-5), which were introduced between matrixes and a ligand.  Among them,
     tartaric acid derivative (5) exhibited the most effective reduction of
nonspecific
     binding proteins, while maintaining binding of the target protein.
     Introduction of 5 on TOYOPEARL reduced tubulin and actin by almost 65% and
     90% compared to that without the hydrophilic spacer, resp., with effective
     binding to the target protein, FKBP12.
IT   675606-56-1P 675606-57-2P 675606-58-3P
     RL: ARU (Analytical role, unclassified); SPN (Synthetic preparation); ANST
     (Analytical study); PREP (Preparation)
        (design and synthesis of novel hydrophilic spacers for reduction of
        nonspecific binding proteins on affinity resins)
RN   675606-56-1  CAPLUS
CN   3,6,9,12,15-Pentaoxaheptadecanoic acid,
     17-[4-[(14-amino-3,6,9,12-tetraoxatetradec-1-yl)oxy]phenoxy]-,
     1,1-dimethylethyl ester  (CA INDEX NAME)
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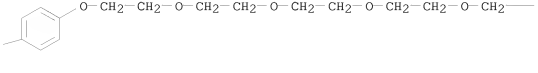
PAGE 1-B



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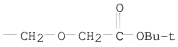

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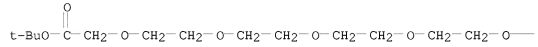

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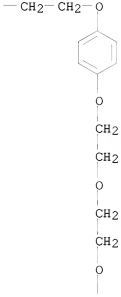

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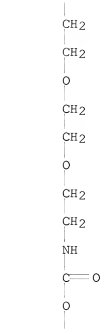

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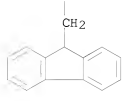

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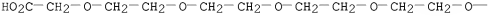
16-[4-[(16-carboxy-3,6,9,12,15-pentaoxahexadec-1-yl)oxy]phenoxy]-,
1-(9H-fluoren-9-ylmethyl) ester (CA INDEX NAME)

PAGE 1-A



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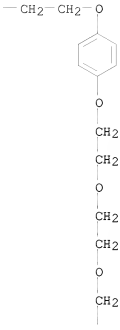

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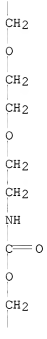

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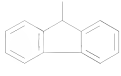

REFERENCE COUNT: 14 THERE ARE 14 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 2 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 2004:252751 CAPLUS
DOCUMENT NUMBER: 140:283950
TITLE: Synthesis of hydrophilic spacers that can reduce nonspecific adsorption of molecules to the surface of solid phase matrix and application to preparation of affinity ligand-immobilized matrixes
INVENTOR(S): Tanaka, Akito; Terada, Tomohiro; Tamura, Tsurunori; Ichiyama, Takaaki; Yamazaki, Akira; Furuya, Minoru; Haramura, Masayuki
PATENT ASSIGNEE(S): Reverse Proteomics Research Institute Co., Ltd., Japan; Fujisawa Pharmaceutical Co., Ltd.
SOURCE: PCT Int. Appl., 134 pp.
CODEN: PIXXD2
DOCUMENT TYPE: Patent
LANGUAGE: Japanese
FAMILY ACC. NUM. COUNT: 1
PATENT INFORMATION:

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2004025297	A1	20040325	WO 2003-JP9640	20030730
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RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
AU 2003254782	A1	20040430	AU 2003-254782	20030730
EP 1553412	A1	20050713	EP 2003-795206	20030730
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
US 20060177943	A1	20060810	US 2006-522716	20060120
PRIORITY APPLN. INFO.:			JP 2002-222226 A	20020730
			WO 2003-JP9640 W	20030730

OTHER SOURCE(S): MARPAT 140:283950

AB Chemical modification method for reducing nonspecific adsorption of mols. to the surface of solid phase matrixes in order to promote specific interaction of immobilized mols. with the target partner mols. The strategy includes the controlling the hydrophobicity of the surface by introducing hydrophilic spacers. These hydrophilic spacers are designed to have ≥ 6 hydrogen bond acceptors, ≥ 5 hydrogen bond donors and their sum are ≥ 9, ≥ 1 carbonyl group, no charged groups. The specific structures for the hydrophilic spacers with reactive carbonyl and amino groups have been designed and synthesized. The hydrophilic spacers have the structures of polyols containing methylenes with linear or branched alkyl group of 1.apprx.3 carbon or -CH2OH, the two or three polyethylene glycol chains (1 .apprx. 1000 EG units) linked to phenol derivs., and alkyl chain (C: n = 1 .apprx. 10) linked with 1 .apprx. 10 unit(s) of (-O-C(R1,R2)C(R3,R4)-) (R1-R4: linear or branched alkyl of 1.apprx.3 carbon). The matrixes with these spacers can be used for immobilization of various sizes of ligands (low or high mol. weight) and for analyses of their interaction with the target partner mols. of various mol. sizes. The applications of some hydrophilic spacers to immobilization of affinity ligands such as FK506 to TOYO-pearl AF-amino resin or gold-film matrixes and detection of FKBP12 protein in rat brain

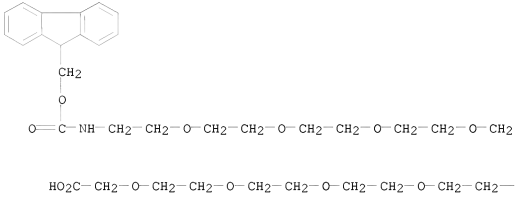
lysate by using the prepared FK506-matrix were demonstrated.

IT 675606-75-4P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(potential use as hydrophilic spacer; synthesis of hydrophilic spacers that can reduce nonspecific adsorption of mols. to surface of solid phase matrix and application to preparation of affinity ligand-immobilized matrixes)

RN 675606-75-4 CAPLUS

CN 5,8,11,14-Tetraoxa-2-azahexadecanoic acid, 16-[3-[(16-carboxy-3,6,9,12,15-pentaoxahexadec-1-yl)oxy]-5-[[18-(9H-fluoren-9-yl)-16-oxo-3,6,9,12,17-pentaoxa-15-azaoctadec-1-yl]oxy]phenoxy]-, 1-(9H-fluoren-9-ylmethyl) ester (CA INDEX NAME)

PAGE 1-A



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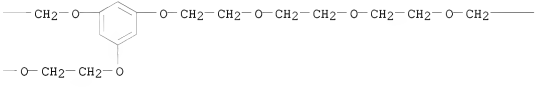

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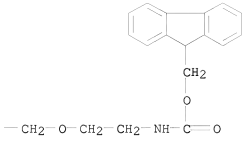

IT 675606-56-1P 675606-57-2P 675606-74-3P
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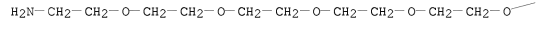
(synthesis of hydrophilic spacers that can reduce nonspecific adsorption of mols. to surface of solid phase matrix and application to preparation of affinity ligand-immobilized matrixes)
RN 675606-56-1 CAPLUS
CN 3,6,9,12,15-Pentaoxaheptadecanoic acid, 17-[4-[(14-amino-3,6,9,12-tetraoxatetradec-1-yl)oxy]phenoxy]-, 1,1-dimethylethyl ester (CA INDEX NAME)

PAGE 1-A



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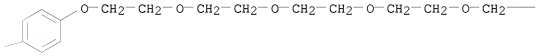

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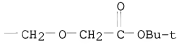
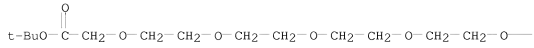


RN 675606-57-2 CAPLUS
CN 3,6,9,12,15-Pentaoxaheptadecanoic acid, 17-[4-[[18-(9H-fluoren-9-yl)-16-oxo-3,6,9,12,17-pentaoxa-15-azaoctadec-1-yl]oxy]phenoxy]-, 1,1-dimethylethyl ester (CA INDEX NAME)

PAGE 1-A



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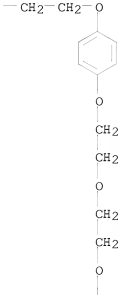

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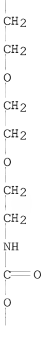

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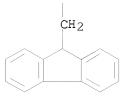

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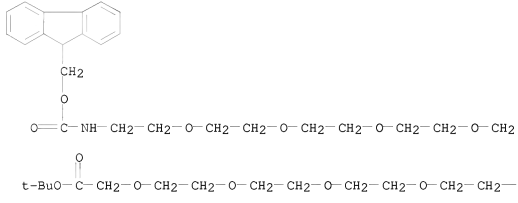
CN 5,8,11,14-Tetraoxa-2-azahexadecanoic acid, 16-[3-[(19,19-dimethyl-17-oxo-3,6,9,12,15,18-hexaoxaeicos-1-yl)oxy]-5-[[18-(9H-fluoren-9-yl)-16-oxo-3,6,9,12,17-pentaoxa-15-azaoctadec-1-yl]oxy]phenoxy]-, 9H-fluoren-9-ylmethyl ester (CA INDEX NAME)

PAGE 1-A



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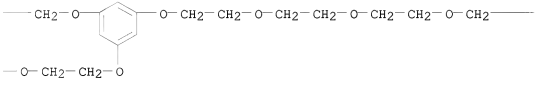

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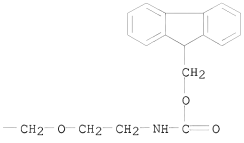
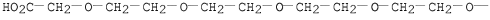


IT 675606-58-3P
RL: BUU (Biological use, unclassified); RCT (Reactant); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)
(used as hydrophilic spacer, linked to TOYO-Pearl resin; synthesis of hydrophilic spacers that can reduce nonspecific adsorption of mols. to surface of solid phase matrix and application to preparation of affinity ligand-immobilized matrixes)
RN 675606-58-3 CAPLUS
CN 5,8,11,14-Tetraoxa-2-azahexadecanoic acid, 16-[4-[(16-carboxy-3,6,9,12,15-pentaoxahexadec-1-yl)oxy]phenoxy]-, 1-(9H-fluoren-9-ylmethyl) ester (CA INDEX NAME)

PAGE 1-A



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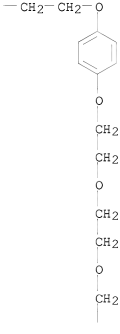
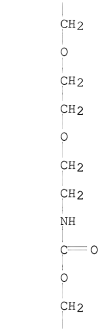
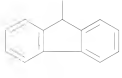






REFERENCE COUNT: 4 THERE ARE 4 CITED REFERENCES AVAILABLE FOR THIS RECORD. ALL CITATIONS AVAILABLE IN THE RE FORMAT

L4 ANSWER 3 OF 3 CAPLUS COPYRIGHT 2009 ACS on STN
ACCESSION NUMBER: 1996:516284 CAPLUS
DOCUMENT NUMBER: 125:247294
ORIGINAL REFERENCE NO.: 125:46221a,46224a
TITLE: Syntheses of ligands containing two and three 2,2'-(bisamino)diphenyl ether units designed for molecular self-assembly on lithiation
AUTHOR(S): Ashton, Peter R.; Hoerner, Bernd; Kocian, Oldrich; Menzer, Stephan; White, Andrew J. P.; Stoddart, J. Fraser; Williams, David J.
CORPORATE SOURCE: School Chem., Univ. Birmingham, Birmingham, B15 2TT, UK
SOURCE: Synthesis (1996), (8), 930-940
CODEN: SYNTBF; ISSN: 0039-7881
PUBLISHER: Thieme
DOCUMENT TYPE: Journal
LANGUAGE: English

AB The syntheses of polyamines containing 2-3 2,2'-(bisamino)diphenyl ether units linked together, designed for self-assembly following lithiation, are reported. The x-ray crystal structures of 2 of the bis[2,2-(bisamino)diphenyl ethers] are described. The ligand, which is linked by an ethylene glycol spacer, exhibits a coiled conformation by intramol. H bonds and supplemented by [CH-π] interactions. The ligand, which is linked by a more rigid bridge, containing a paraphenylene unit,

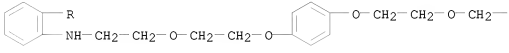
displays a stretched conformation stabilized by intramol. edge to face interactions.

IT 181725-62-2P
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)
(preparation of ligands with (bisamino)diphenyl ether units)

RN 181725-62-2 CAPLUS

CN Acetamide, 2,2'-[oxybis(2,1-phenyleneimino-2,1-ethanediyloxy-2,1-ethanediyloxy-4,1-phenyleneoxy-2,1-ethanediyloxy)]bis[N-[2-[2-[(methoxyacetyl)amino]phenoxy]phenyl]- (9CI) (CA INDEX NAME)

PAGE 1-A



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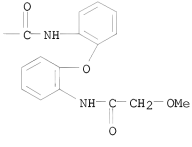
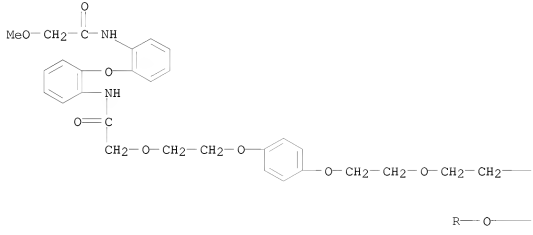


PAGE 2-A




NH